CLAIM LISTING

1. (Previously presented) A compound of formula (Ia)

$$R^{2}$$
 R^{3}
 R^{4}
 $(CH_{2})_{n}$
 R^{5}
 $(O)_{m}$
 R^{6}
 $(OR^{7})_{n}$
 (Ia)

wherein R^1 , R^2 , R^3 , and R^4 independently of each other represent hydrogen, halogen, perhalomethyl, hydroxy, nitro, cyano, formyl, or C_{1-12} -alkyl, C_{4-12} -alkenynyl, C_{2-12} -alkenyl, C_{2-12} -alkynyl, C_{1-12} -alkoxy, aryl, aryloxy, aralkyl, aralkoxy, heterocyclyl, heteroaryl, heteroaralkyl, heteroaryloxy, heteroaralkoxy, acyl, acyloxy, hydroxy C_{1-12} -alkyl, amino, acylamino, C_{1-12} -alkylamino, arylamino, aralkylamino, amino C_{1-12} -alkyl, C_{1-12} -alkoxycarbonyl, aryloxycarbonyl, aralkoxycarbonyl, C_{1-12} -alkoxy C_{1-12} -alkyl, aryloxy C_{1-12} -alkyl, aralkoxy C_{1-12} -alkyl, C_{1-12} -alkyl, C_{1-12} -alkyl, C_{1-12} -alkyl, C_{1-12} -alkyl, C_{1-12} -alkyl, C_{1-12} -alkoxycarbonylamino, aryloxycarbonylamino, aralkoxycarbonylamino, - COR^{11} , or - SO_2R^{12} , wherein R^{11} and R^{12} independently of each other are selected from hydroxy, halogen, perhalomethyl, C_{1-6} -alkoxy or amino optionally substituted with one or more C_{1-6} -alkyl, perhalomethyl or aryl; optionally substituted with one or more halogen, perhalomethyl, hydroxy, nitro or cyano; or R^1 and R^2 , R^2 and R^3 and/or R^3 and R^4 may form a cyclic ring containing from 5 to

or R^1 and R^2 , R^2 and R^3 and/or R^3 and R^4 may form a cyclic ring containing from 5 to 7 carbon atoms optionally substituted with one or more C_{1-6} -alkyl;

ring A fused to the ring containing X and N represents a 5-6 membered cyclic ring, optionally substituted with one or more hydrogen, halogen, perhalomethyl, hydroxy or C₁₋₇-alkyl, C₂₋₇-alkenyl, C₂₋₇-alkynyl, C₁₋₇-alkoxy or aryl;

X is -O-(CHR⁹)-, -O-CH₂-O-, -CH₂-O-CH₂-, wherein R⁹ is hydrogen, halogen, hydroxy, nitro, cyano, formyl, C₁₋₁₂-alkyl, C₁₋₁₂-alkoxy, aryl, aryloxy, aralkyl, aralkoxy, heterocyclyl, heteroaryl, heteroaralkyl, heteroaryloxy, heteroaralkoxy, acyl, acyloxy, hydroxyalkyl, amino, acylamino, C₁₋₁₂-alkylamino, arylamino, aralkylamino, aminoC₁₋₁₂-alkyl, C₁₋₁₂-alkoxycarbonyl, aryloxycarbonyl, aralkoxycarbonyl, C₁₋₁₂-alkoxyC₁₋₁₂-alkyl, aryloxyC₁₋₁₂-alkyl, aralkoxyC₁₋₁₂-alkyl, C₁₋₁₂-alkylthio, thioC₁₋₁₂-alkyl, C₁₋₁₂-alkoxycarbonylamino, aryloxycarbonylamino, aralkoxycarbonylamino, -COR¹³, or -SO₂R¹⁴, wherein R¹³ and R¹⁴ independently of each other are selected from hydroxy, halogen, C₁₋₆-alkoxy, amino optionally substituted with one or more C₁₋₆-alkyl, perhalomethyl or aryl;

Ar represents arylene or heteroarylene, optionally substituted with one or more C₁₋₆-alkyl or aryl;

 R^5 represents hydrogen, hydroxy, halogen, C_{1-12} -alkoxy, C_{1-12} -alkyl, C_{4-12} -alkenynyl, C_{2-12} -alkenyl, C_{2-12} -alkynyl or aralkyl; optionally substituted with one or more halogen, perhalomethyl, hydroxy, nitro or cyano; or R^5 forms a bond together with R^6 , R^6 represents hydrogen, hydroxy, halogen, C_{1-12} -alkoxy, C_{1-12} -alkyl, C_{4-12} -alkenynyl, C_{2-12} -alkenyl, C_{2-12} -alkynyl, acyl or aralkyl; optionally substituted with one or more halogen, perhalomethyl, hydroxy, nitro or cyano; or R^6 forms a bond together with R^5 , R^7 represents hydrogen, C_{1-12} -alkyl, C_{4-12} -alkenynyl, C_{2-12} -alkenyl, C_{2-12} -alkynyl, aryl, aralkyl, C_{1-12} -alkoxy C_{1-12} -alkoxycarbonyl, aryloxycarbonyl, C_{1-12} -alkylaminocarbonyl, arylaminocarbonyl, acyl, heterocyclyl, heteroaryl or heteroaralkyl groups; optionally substituted with one or more halogen, perhalomethyl, hydroxy, nitro or cyano; R^8 represents hydrogen, C_{1-12} -alkyl, C_{4-12} -alkenynyl, C_{2-12} -alkenyl, C_{2-12} -alkynyl, aryl, aralkyl, heterocyclyl, heteroaryl or heteroaralkyl groups; optionally substituted with one or more halogen, perhalomethyl, hydroxy, nitro or cyano;

Y represents oxygen, sulphur or NR^{10} , where R^{10} represents hydrogen, C_{1-12} -alkyl, aryl, hydroxy C_{1-12} -alkyl or aralkyl groups or when Y is NR^{10} , R^8 and R^{10} may form a 5 or 6 membered nitrogen containing ring, optionally substituted with one or more C_{1-6} -alkyl; n is an integer ranging from 1 to 4 and m is an integer ranging from 0 to 1; or a pharmaceutically acceptable salt thereof.

- 2. (Original) A compound according to claim 1 wherein R^1 , R^2 , R^3 , and R^4 independently of each other represent hydrogen, halogen, perhalomethyl, hydroxy, cyano, or C_{1-7} -alkyl, C_{4-7} -alkenynyl, C_{2-7} -alkenyl, C_{2-7} -alkynyl, C_{1-7} -alkoxy, aryl, aryloxy, aralkyl, aralkoxy, heterocyclyl, heteroaryl, heteroaralkyl, heteroaryloxy, heteroaralkoxy, acyl, acyloxy, hydroxy C_{1-7} -alkyl, amino, acylamino, C_{1-7} -alkylamino, arylamino, aralkylamino, amino C_{1-7} -alkyl, C_{1-7} -alkoxy C_{1-7} -alkyl, aryloxy C_{1-7} -alkyl, aralkoxy C_{1-7} -alkyl, C_{1-7} -alkyl, C_{1-7} -alkyl, C_{1-7} -alkyl, C_{1-7} -alkoxycarbonylamino, aralkoxycarbonylamino, -COR 11 , or -SO $_2$ R 12 , wherein R 11 and R 12 independently of each other are selected from hydroxy, perhalomethyl or amino optionally substituted with one or more C_{1-6} -alkyl, perhalomethyl or aryl; optionally substituted with one or more halogen, perhalomethyl, hydroxy or cyano; or R 1 and R 2 , R 2 and R 3 and/or R 3 and R 4 may form a cyclic ring containing from 5 to 7 carbon atoms optionally substituted with one or more C_{1-6} -alkyl.
- 3. (Cancelled)
- 4. (Cancelled)
- 5. (Cancelled)
- 6. (Cancelled)
- 7. (Previously presented) A compound according to claim 1 wherein ring A fused to the ring containing X and N represents a 5-6 membered cyclic ring, optionally substituted with one or more hydrogen, halogen, perhalomethyl, hydroxy, cyano, or C₁₋₇-alkyl, C₄₋₇-alkenynyl, C₂₋₇-alkenyl, C₂₋₇-alkynyl, C₁₋₇-alkoxy, aryl, aryloxy, aralkyl, aralkoxy, heterocyclyl, heteroaryl, heteroaralkyl, heteroaryloxy, heteroaralkoxy, acyl, acyloxy, hydroxyC₁₋₇-alkyl, amino, acylamino, C₁₋₇-alkylamino, arylamino, aralkylamino, aminoC₁₋₇-alkyl, C₁₋₇-alkoxyC₁. 7-alkyl, aryloxyC₁₋₇-alkyl, aralkoxyC₁₋₇-alkyl, C₁₋₇-alkylthio, thioC₁₋₇-alkyl, C₁₋₇-alkoxyC₁-alkoxyCarbonylamino, aryloxycarbonylamino, aralkoxycarbonylamino, -COR¹¹, or -SO₂R¹², wherein R¹¹ and R¹² independently of each other are selected from hydroxy, perhalomethyl or amino optionally substituted with one or more C₁₋₆-alkyl, perhalomethyl or aryl; optionally substituted with one or more halogen, perhalomethyl, hydroxy or cyano.

- 8. (Cancelled)
 9. (Cancelled)
 10. (Cancelled)
 11. (Cancelled)
 12. (Cancelled)
 13. (Cancelled)
- 14. (Cancelled)
- 15. (Cancelled)
- 16. (Cancelled)
- 17. (Previously presented) A compound according to claim 1 wherein Ar represents arylene or heteroarylene;

R⁵ represents hydrogen, hydroxy, halogen; or R⁵ forms a bond together with R⁶,

R⁶ represents hydrogen, hydroxy, halogen; or R⁶ forms a bond together with R⁵,

 R^7 represents hydrogen, C_{1-7} -alkyl, C_{2-7} -alkenyl, C_{2-7} -alkynyl, aryl, aralkyl, C_{1-7} -alkoxy C_{1-7} -alkyl, C_{1-7} -alkylaminocarbonyl, arylaminocarbonyl, acyl, heterocyclyl, heteroaryl or heteroaralkyl groups;

R⁸ represents hydrogen, C₁₋₇-alkyl, C₂₋₇-alkenyl, C₂₋₇-alkynyl;

Y represents oxygen or sulphur;

n is an integer ranging from 2 to 3 and m is 1.

- 18. (Cancelled)
- 19. (Cancelled)
- 20. (Cancelled)
- 21. (Cancelled)
- 22. (Cancelled)
- 23. (Cancelled)
- 24. (Cancelled)
- 25. (Cancelled)
- 26. (Cancelled)
- 27. (Cancelled)

- 28. (Cancelled)
- 29. (Cancelled)
- 30. (Cancelled)
- 31. (Cancelled)
- 32. (Cancelled)
- 33. (Cancelled)
- 34. (Cancelled)
- 35. (Cancelled)
- 36. (Cancelled)
- 37. (Cancelled)
- 38. (Cancelled)
- 39. (Cancelled)
- 40. (Cancelled)
- 41. (Cancelled)
- 42. (Cancelled)
- 43. (Cancelled)
- 44. (Cancelled)
- 45. (Previously presented) The compound according to claim 1 which is
- 2-Ethoxy-3-(4-[2-(5,11-dihydro-5H-dibenzo[b,e][1,4]oxazepin-5-yl)-ethoxy]-phenyl)-propionic acid,
- 2-Methoxy-3-(4-[2-(5,11-dihydro-5H-dibenzo[b,e][1,4]oxazepin-5-yl)-ethoxy]-phenyl)-propionic acid,
- 2-Propoxy-3-(4-[2-(5,11-dihydro-5H-dibenzo[b,e][1,4]oxazepin-5-yl)-ethoxy]-phenyl)-propionic acid,
- 2-Benzyloxy-3-(4-[2-(5,11-dihydro-5H-dibenzo[b,e][1,4]oxazepin-5-yl)-ethoxy]-phenyl)-propionic acid,
- 2-Ethoxy-3-(4-[3-(5,11-dihydro-5H-dibenzo[b,e][1,4]oxazepin-5-yl)-propionic acid,
- 2-Methoxy-3-(4-[3-(5,11-dihydro-5H-dibenzo[b,e][1,4]oxazepin-5-yl)-propoxy]-phenyl)-propionic acid,

- 2-Benzyloxy-3-(4-[3-(5,11-dihydro-5H-dibenzo[b,e][1,4]oxazepin-5-yl)-propoxy]-phenyl)-propionic acid,
- 2-Ethoxy-3-(4-[3-(5,11-dihydro-5H-dibenzo[b,e][1,4]oxazepin-5-yl)-propyl]-phenyl)-propionic acid,
- 2-Methoxy-3-(4-[3-(5,11-dihydro-5H-dibenzo[b,e][1,4]oxazepin-5-yl)-propyl]-phenyl)-propionic acid,
- 2-Benzyloxy-3-(4-[3-(5,11-dihydro-5H-dibenzo[b,e][1,4]oxazepin-5-yl)-propyl]-phenyl)-propionic acid,
- 2-Ethoxy-3-(4-[1-(5,11-dihydro-5H-dibenzo[b,e][1,4]oxazepin-5-yl)-methoxy]-phenyl)-propionic acid,
- $3-\{4-[2-(6,7-Dihydro-5H-dibenzo[b,g]azocin-12-yl)-ethoxy]-phenyl\}-2-ethoxy-propionic acid,$
- $3-\{4-[2-(6,7-Dihydro-5H-dibenzo[b,g]azocin-12-yl)-ethoxy]-phenyl\}-2-propoxy-propionic acid,$
- 3-{4-[2-(6,7-Dihydro-5*H*-dibenzo[*b*,*g*]azocin-12-yl)-ethoxy]-phenyl}-2-methoxy-propionic acid,
- 3-{4-[2-(6,7-Dihydro-5*H*-dibenzo[*b*,*g*]azocin-12-yl)-ethoxy]-phenyl}-2-benzyloxy-propionic acid,
- $3-\{4-[1-(6,7-\text{Dihydro}-5H-\text{dibenzo}[b,g]\text{azocin}-12-yl)-\text{methoxy}]-\text{phenyl}\}-2-\text{ethoxy-propionic}$ acid,
- $3-\{4-[3-(6,7-Dihydro-5H-dibenzo[b,g]azocin-12-yl)-propoxy]-phenyl\}-2-ethoxy-propionic acid,$
- 3-{4-[3-(6,7-Dihydro-5*H*-dibenzo[*b*,*g*]azocin-12-yl)-propoxy]-phenyl}-2-methoxy-propionic acid,
- $3-\{4-[3-(6,7-Dihydro-5H-dibenzo[b,g]azocin-12-yl)-propoxy]-phenyl\}-2-benzyloxy-propionic acid,$
- $3-\{4-[3-(6,7-Dihydro-5H-dibenzo[b,g]azocin-12-yl)-propyl]-phenyl\}-2-ethoxy-propionic acid,$
- $3-\{4-[3-(6,7-Dihydro-5H-dibenzo[b,g]azocin-12-yl)-propyl]-phenyl\}-2-methoxy-propionic acid,$

- $3-\{4-[3-(6,7-Dihydro-5H-dibenzo[b,g]azocin-12-yl)-propyl]-phenyl\}-2-benzyloxy-propionic acid,$
- 3-(4-Dibenzo[d,g]dioxazocin-12-yl)-1-propoxy)-phenyl-2-ethoxy-propionic acid,
- 3-(4-Dibenzo[d,g]dioxazocin-12-yl)-1-propoxy)-phenyl-2-methoxy-propionic acid,
- 3-(4-Dibenzo[d,g]dioxazocin-12-yl)-1-propoxy)-phenyl-2-propoxy-propionic acid,
- 3-(4-Dibenzo[d,g]dioxazocin-12-yl)-1-propoxy)-phenyl-2-benzyloxy-propionic acid,
- 3-(4-Dibenzo[d,g]dioxazocin-12-yl)-1-propyl)-phenyl-2-ethoxy-propionic acid,
- 3-(4-Dibenzo[d,g]dioxazocin-12-yl)-1-propyl)-phenyl-2-methoxy-propionic acid,
- 3-(4-Dibenzo[d,g]dioxazocin-12-yl)-1-propyl)-phenyl-2-propoxy-propionic acid,
- 3-(4-Dibenzo[d,g]dioxazocin-12-yl)-1-propyl)-phenyl-2-benzyloxy-propionic acid,
- 2-(4-Dibenzo[d,g]dioxazocin-12-yl)-1-ethoxy)-phenyl-2-ethoxy-propionic acid,
- 2-(4-Dibenzo[d,g]dioxazocin-12-yl)-1-ethoxy)-phenyl-2-propoxy-propionic acid,
- 1-(4-Dibenzo[d,g]dioxazocin-12-yl)-1-methoxy)-phenyl-2-ethoxy-propionic acid,
- 2-(4-Dibenzo[d,g]dioxazocin-12-yl)-1-ethoxy)-phenyl-2-benzyloxy-propionic acid, or a pharmaceutically acceptable salt thereof.
- 46. (Previously presented) The compound according to claim 1 which is
- 2-Ethoxy-3-{4-[2-(5,11-dihydro-5H-dibenzo[b,e][1,4]oxazepin-5-yl)-ethoxy]-phenyl}-propionic acid,
- $3-\{4-[2-(6,7-Dihydro-5H-dibenzo[b,g]azocin-12-yl)-ethoxy]-phenyl\}-2-ethoxy-propionic acid,$
- or a pharmaceutically acceptable salt thereof.
- 47. (Previously presented) A pharmaceutical composition comprising, as an active ingredient, a compound according to claim 1 or a pharmaceutically acceptable salt thereof together with a pharmaceutically acceptable carrier or diluent.
- 48. (Cancelled)
- 49. (Cancelled)
- 52. (Cancelled)
- 51. (Cancelled)
- 52. (Cancelled)

53. (Cancelled)

- 54. (Currently amended) A method for the treatment of conditions mediated by the Peroxisome Proliferator-Activated Receptors (PPAR), wherein the condition is selected from the following: type 2 diabetes, impaired glucose tolerance, hypertension, obesity, insulin resistance, hyperglycemia, atherosclerosis, hyperlipidemia, coronary artery disease, glomerulonephritis, glomerulosclerosis, nephritic syndrome, hypertensive nephrosclerosis, dementia, diabetic complications, psoriasis, polycystic ovarian syndrome and osteoporosis, the method comprising administering to a subject in need thereof an effective amount of a compound according to claim 1 or a pharmaceutically acceptable salt thereof.
- 55. (Previously presented) A method for the treatment of diabetes or obesity, the method comprising administering to a subject in need thereof an effective amount of a compound according to claim 1 or a pharmaceutically acceptable salt thereof.
- 56. (Cancelled)
- 57. (Cancelled)
- 58. (Cancelled)
- 59. (Cancelled)
- 60. (Cancelled)